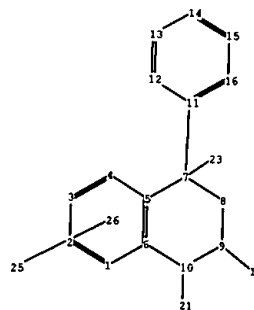


c # 1

L15



100 1

chain nodes :

21 23

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16

ring/chain nodes :

17 18 25

chain bonds :

7-11 7-23 9-17 10-21

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-10 7-8 8-9 9-10 11-12 11-16
12-13 13-14 14-15 15-16

exact/norm bonds :

5-7 6-10 7-8 7-23 8-9 9-10 9-17 10-21

exact bonds :

7-11

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6 11-12 11-16 12-13 13-14 14-15 15-16

isolated ring systems :

containing 1 :

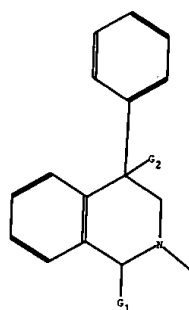
G1:H, [*1]

G2:H,X

Match level :

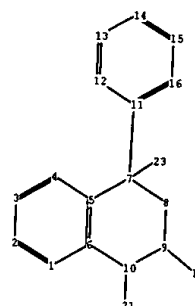
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom
10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:CLASS
18:CLASS

21:CLASS 24:CLASS 25:CLASS 26:CLASS



c 0¹

L4



100¹

chain nodes :

21 23

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16

ring/chain nodes :

17 18

chain bonds :

7-11 7-23 9-17 10-21

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-10 7-8 8-9 9-10 11-12 11-16
12-13 13-14 14-15 15-16

exact/norm bonds :

5-7 6-10 7-8 7-23 8-9 9-10 9-17 10-21

exact bonds :

7-11

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6 11-12 11-16 12-13 13-14 14-15 15-16

G1:H, [*1]

G2:H, X

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom
10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:CLASS
18:CLASS 21:CLASS 23:CLASS